IN THE CLAIMS

The following is an updated listing of the claims in the application with claims 1, 2, 3, 5, 10 and 12 shown as currently amended and claim 13 as cancelled:

LISTING OF CLAIMS

1. (Currently amended) An indene derivative of formula (I) or a pharmaceutically acceptable salt thereof:

wherein,

 R_1 is C_{1-6} alkyl, C_{1-6} alkenyl, or C_{3-6} cycloalkyl, <u>each of</u> which is unsubstituted or substitutied with one or more phenyl groups;

$$R_2$$
 is H, CN, CO_2R^a , $CH_2CO_2R^a$, $CONR^bR^c$, R_3 is C_{1-6} alkyl, C_{3-6} cycloalkyl, $\frac{\partial^2}{\partial r^2}$ naphthyl, phenyl,

unsubstituted or substitutied with one or more substituents selected from the group consisting of halogen, CN, NH₂, NO₂, OR^a, phenyloxy, C₁₋₆ alkyl, and C₃₋₆ cycloalkyl; and

R₄, R₅, R₆, and R₇ are each independently H, OH, OSO₂CH₃, O(CH₂)_mR^e, CH₂R^f, OCOCH₂OR^g, OCH₂CH₂OR^g of, OCH₂CH=CHR^g, or pyridine-2-yloxy, or R₅ and R₆ together form OCH₂O;

in which R^a is H, or C_{1-6} alkyl, or C_{3-6} cycloalkyl, which is C_{1-6} alkyl and C_{3-6} cycloalkyl being each unsubstituted or substituted with one or more halogens;

R^b and R^c are each independently H, C₁₋₆ alkyl, or C₃₋₆ cycloalkyl;

R^d is O, S, or NR^a;

R^e is H, halogen, C₃₋₆ cycloalkyl, naphthyl,

$$-\frac{1}{2} \left\{ \begin{array}{c} R^{a} \\ N \end{array} \right\} = \left\{ \begin{array}{ccc} R^{d} \\ \end{array} \right\} \left\{ \begin{array}{ccc} R^{d} \\ \end{array} \right\} \left\{ \begin{array}{ccc} R^{a} \\ \end{array} \right\} \left\{ \begin{array}{ccc}$$

<u>, adamantly</u>, or phenyl, which

is phenyl being unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, CN, NH₂, NO₂, OR^a, CF₃, and COOR^a;

$$R^f$$
 is $OCH_2CH_2R^g$ or $-\frac{1}{2}$ -N R^d ;

R^g is phenyl, which is unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, CN, NH₂, NO₂, and OR^a; and

m is an integer in the range of 1 to 5.

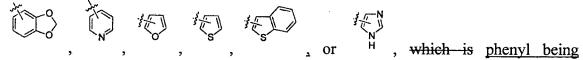
2. (Currently amended) The compound of claim 1, wherein R_1 is C_{1-6} alkyl, which is unsubstituted or substituted with a phenyl group; R_2 is H, CN, CO_2R^a , $CH_2CO_2R^a$, $CONR^bR^c$, or phenyl; R_3 is C_{1-6} alkyl, C_{3-6} cycloalkyl, Θ phenyl,

unsubstituted or substitutied with one or more substituents selected from the group consisting of halogen, C_{1-6} alkyl, and C_{3-6} cycloalkyl; R_4 and R_7 are H; R_5 and R_6 are each independently OH, OSO_2CH_3 , $O(CH_2)_mR^e$, CH_2R^f , $OCOCH_2OR^g$, $OCH_2CH_2OR^g$, or $OCH_2CH=CHR^g$, or together form OCH_2O ; R^a is H_7 or C_{1-6} alkyl; R^d is O or NCH_3 ; R^e is H, halogen, C_{3-6} cycloalkyl, naphthyl,

being unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, OH, methoxy, CF₃, and COOR^a; R^f is OCH₂CH₂R^g or

$$-$$
{-N R^d ; and R^g is phenyl.

3. (Currently amended) The compound of claim 2, wherein R_1 is CH_3 ; R_2 is H_3 , CN_4 , CO_2R^a , or $CONR^bR^c$; R_3 is C_{1-6} alkyl, Θ_F phenyl,



unsubstituted or substitutied with one or more halogens or C_{1-6} alkyl groups; and R_5 and R_6 are each independently $O(CH_2)_m R^e$ or $CH_2 R^f$, or together form $OCH_2 O$.

- 4. (As originally filed) The compound of claim 1, which is selected from the group consisting of:
- 1) 6-methoxy-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 2) 1-(*trans*-isopropylimino-*N*-oxy)-6-methoxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 3) 1-(*trans*-benzylimino-*N*-oxy)-6-methoxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 4) 1-(*trans*-ethylimino-*N*-oxy)-6-methoxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 5) 6-methoxy-1-(*trans*-phenylpropylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 6) 6-methoxy-1-(*trans*-(2-methylbutenylimino)-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 7) 1-(*trans*-isobutylimino-*N*-oxy)-6-methoxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 8) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morphorline-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 9) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 10) 1-(*trans*-methylimino-*N*-oxy)-6-phenetyloxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 3-furan-3-yl-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 12) 6-hydroxy-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 13) 1-(*cis*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 14) 3-(trans-methylimino-N-oxy)-1-phenyl-3H-indene-5-ol
- 15) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(5-phenylpentyloxy)-1H-indene-2-carboxylate ethyl ester
- 16) 1-(*cis*-methylimino-*N*-oxy)-3-phenyl-6-(5-phenylpentyloxy)-1H-indene-2-carboxylate ethyl ester
- 17) 6-[2-(4-chlorophenoxy)acetoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester

18) 6-[2-(4-chlorophenoxy)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester

- 19) 1-(*trans*-methylimino-*N*-oxy)-6-(naphthalene-2-ylmethoxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 20) methyl-[3-phenyl-6-(3-phenylpropoxy)indene-1-yllidene]amine-N-oxide
- 21) 1-(*trans*-methylimino-*N*-oxy)-6-[2-(5-methyl-2-phenylthiazol-4-yl)ethoxy]-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 22) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 23) 6-[2-(4-hydroxyphenyl)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 24) __6-(2-adaman-1-ylethoxy)-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 25) 6-(2-cyclohexylethoxy)-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 26) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylprophenoxy)-1H-indene-2-carboxylate ethyl ester
- 27) 6-[2-(2-fluorophenyl)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 28) 6-[2-(3-fluorophenyl)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 29) 6-[2-(4-fluorophenyl)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 30) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-[2-(3-trifluoromethylphenyl)ethoxy]-1H-indene-2-carboxylate ethyl ester
- 31) 6-(4-methoxycarbonylbenzyloxy)-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 32) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl amide
- 33) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 34) 6-[2-(cyclohexylmethylamino)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 35) 3-(2-fluorophenyl)-6-methoxy-1-(*trans*-methylimino-*N*-oxy)-1H-indene-2-carboxylate ethyl ester
- 36) 1-(*trans*-methylimino-*N*-oxy)-6-[2-(4-methylpiperazine-1-yl)ethoxy]-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 37) (2,3-diphenyl indene-1-yl lidene)methylamine-N-oxide
- 38) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate isopropyl amide
- 39) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate cyclohexyl amide
- 40) [1-(trans-methylimino-N-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-

yl]morpholine-4-yl-methanone

41) 1-(trans-methylimino-N-oxy)-6-(2-morpholine-4-yl-ethoxy)-3-phenyl-1H-indene-2-carboxylate cyclohexyl amide

- 42) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-5-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 43) 1-(*trans*-methylimino-*N*-oxy)-6-phenethyloxymethyl-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 44) (6-methoxy-3-phenylindene-1-yllidene)methylamine-N-oxide
- 45) 1-(cis-methylimino-N-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 6-(2-bromoethoxy)-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 47) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate *tert*-buthyl ester
- 48) 1-(*trans*-methylimino-*N*-oxy)-5,6-methylenedioxy-1-oxo-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 49) 4-[2-isopropylcarbamoyl-3-(*trans*-methylimino-*N*-oxy)-1-phenyl-3H-indene-5-yl-oxylmethyl]benzoate methyl ester
- 50) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate isopropyl amide
- 51) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate cyclopropyl amide
- 52) 3-(3-fluorophenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 53) (6-methoxy-1-(trans-methylimino-N-oxy)-3-phenyl-1H-indene-2-yl)acetate ethyl ester
- 54) (6-methoxy-1-(cis-methylimino-N-oxy)-3-phenyl-1H-indene-2-yl)acetate ethyl ester
- 55) 5-[2-(5-ethylpyridine-2-yl)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate isopropyl amide
- 56) 1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-3-*p*-tolyl-1H-indene-2-carboxylate ethyl ester
- 57) 1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-3-thiophene-2-yl-1H-indene-2-carboxylate ethyl ester
- 58) 3-(4-chlorophenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 59) 3-(5-chlorothiophene-2-yl)-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenyl propoxy)-1H-indene-2-carboxylate ethyl ester
- 60) 1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-3-*m*-tolyl-1H-indene-2-carboxylate ethyl ester
- 61) 1-(*trans*-methylimino-*N*-oxy)-3-(4-phenoxyphenyl)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 62) 3-benzo-[1,3]-dioxol-5-yl-1-(trans-methylimino-N-oxy)-6-(3-phenyl propoxy)-

- 1H-indene-2-carboxylate ethyl ester
- 63) methyl-[6-(3-phenylpropoxy)-3-pyridine-2-yl-indene-1-yllidene]-amine-N-oxide
- 3-furan-2-yl-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 65) 3-ethyl-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 3-methyl-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 67) 1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-3-thiophene-3-yl-1H-indene-2-carboxylate ethyl ester
- 68) 3-cyclopropyl-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 69) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-thiophene-3-yl-1H-indene-2-carboxylate ethyl ester
- 70) 3-benzo-[b]-thiophene-3-yl-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenyl propoxy)-1H-indene-2-carboxylate ethyl ester
- 71) 3-(1H-imidazole-4-yl)-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 72) 3-(1-ethyl propyl)-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 73) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate amide
- 74) 6-(4-benzylmorpholine-2-ylmethoxy)-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate isopropyl amide
- 75) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carbonitrile
- 76) 1-(*trans*-methylimino-*N*-oxy)-5,6-methylenedioxy-1-oxo-3-phenyl-1H-phenyl-2-carboxylate isopropyl amide
- 77) 1-(*trans*-methylimino-*N*-oxy)-6-morpholine-4-ylmethyl-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 78) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate ethyl ester
- 79) 6-[2-(5-ethylpyridine-2-yl)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 80) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 81) 6-[2-(5-ethylpyridine-2-yl)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate isopropyl amide
- 82) methyl-[6-(2-morpholine-4-ylethoxy)-3-phenylindene-1-yllidene]amine-N-oxide
- 83) 5,6-bis-methanesulfonyloxy-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 84) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate isobutyl ester

85) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate methyl ester

- 86) 1-(*cis*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate methyl ester
- 87) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate propyl ester
- 88) 3-(4-fluorophenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate ethyl ester
- 89) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(pyridine-2-ylmethoxy)-1H-indene-2-carboxylate ethyl ester
- 90) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(pyridine-2-yloxy)-1H-indene-2-carboxylate ethyl ester
- 91) 6-(3-methoxybenzyloxy)-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 92) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-thiophene-3-yl-1H-indene-2-carboxylate isopropyl amide
- 93) 3-(1-ethylpropyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-yl ethoxy)-1H-indene-2-carboxylate ethyl ester
- 3-benzo-[b]-thiophene-3-yl-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 95) 3-(4-fluorophenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 96) 3-(1-ethylpropyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 97) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-(2,4,6-trimethylphenyl)-1H-indene-2-carboxylate ethyl ester
- 98) 3-(2,6-dimethylphenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate ethyl ester
- 99) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-5-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 100) 1-(*trans*-methylimino-*N*-oxy)-5-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate isopropyl amide
- 101) 1-(*cis*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate isopropyl ester
- 102) 3-(3-fluorophenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 103) 6-[2-(5-ethylpyridine-2-yl)ethoxy]-3-(3-fluorophenyl)-1-(*trans*-methylimino-*N*-oxy)-1H-indene-2-carboxylate isopropyl amide
- 104) 3-(4-cyanophenyl)-6-(2-morpholine-4-ylethoxy)-1-(*trans*-methylimino-*N*-oxy)-1H-indene-2-carboxylate ethyl ester
- 105) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate isopropyl ester.

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5. (Currently amended) A process for preparing the indene derivative of claim 1, which comprises the step of subjecting an indenone compound of formula (II) to a condensation reaction with R₁NHOH to obtain a compound of formula (II); or comprises the steps of subjecting an indenone compound of formula (III) to a condensation reaction with NH₂OH to obtain a compound of formula (III), and conducting a reaction of the compound of formula (III) with R₁X to obtain a compound of formula (I):

$$\begin{array}{c} R_{q} & O_{q} \\ R_{q} & R_{q} \\ R_{q} & R_{q} \end{array}$$

$$(I)$$

wherein,

X is halogen;

 R_1 is C_{1-6} alkyl, C_{1-6} alkenyl, or C_{3-6} cycloalkyl, <u>each of</u> which is unsubstituted or substitutied with one or more phenyl groups;

, which is phenyl and

unsubstituted or substitutied with one or more substituents selected from the group consisting of halogen, CN, NH₂, NO₂, OR^a, phenyloxy, C₁₋₆ alkyl, and C₃₋₆ cycloalkyl; and

R₄, R₅, R₆₂ and R₇ are each independently H, OH, OSO₂CH₃, O(CH₂)_mR^e, CH₂R^f, OCOCH₂OR^g, OCH₂CH₂OR^g of, OCH₂CH=CHR^g, or pyridine-2-yloxy, or R₅ and R₆ together form OCH₂O;

in which Ra is H, C1-6 alkyl, or C3-6 cycloalkyl, which is C1-6 alkyl and C3-6 cycloalkyl being each unsubstituted or substitutied with one or more halogens;

 R^b and R^c are each independently H, C_{1-6} alkyl, or C_{3-6} cycloalkyl;

R^d is O, S, or NR^a; R^e is H, halogen, C₃₋₆ cycloalkyl, naphthyl,

 $-\frac{1}{2} \underbrace{ \left(\begin{array}{ccc} R^a \\ \end{array} \right)^{R^a}}_{N} - \underbrace{ \left(\begin{array}{ccc} R^d \\ \end{array} \right)^{R^a}}_{N} \underbrace{ \left(\begin{array}{ccc} R^d \\ \end{array} \right)^{R^a}}_{R^a} \underbrace{ \left(\begin{array}{ccc} R^d \\ \end{array} \right)^{R^a}}_{N} \underbrace$

, adamantly, or phenyl, which is phenyl being unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, CN, NH₂, NO₂, OR^a, CF₃, and COOR^a;

$$R^f \text{ is } OCH_2CH_2R^g \text{ or } \overset{-\frac{1}{2}-N}{ } \overset{R^d}{\longrightarrow} ;$$

Rg is phenyl, which is unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, CN, NH2, NO2, and ORa; and

m is an integer in the range of 1 to 5.

- 6. (As originally filed) The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:
 - 1) reacting compounds of formula (V) and (VI) to obtain a compound of formula (VII);
 - 2) subjecting the compound of formula (VII) to cyclization to obtain a compound of formula (VIII); and
 - 3) subjecting the compound of formula (VIII) to oxidation.

$$R_{3}$$
 R_{2}
 R_{5}
 R_{4}
 R_{5}
 R_{5}
 R_{4}
 R_{5}
 R_{5}
 R_{6}
 R_{7}
 R_{8}
 R_{8}
 R_{8}
 R_{8}
 R_{8}
 R_{8}
 R_{8}
 R_{8}
 R_{8}
 R_{9}
 R_{9}
 R_{1}
 R_{2}
 R_{2}
 R_{3}
 R_{4}
 R_{2}
 R_{3}
 R_{4}
 R_{5}
 R_{4}
 R_{5}
 R_{5}
 R_{4}
 R_{5}
 R_{5}

wherein,

 R_2 to R_7 have the same meanings as defined in claim 5, and Z is halogen or activated leaving group.

- 7. (As originally filed) The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:
 - 1) reacting compounds of formula (IX) and (X) to obtain a compound of formula (XI);
 - 2) subjecting the compound of formula (XI) to cyclization to obtain a compound of formula (XII); and
 - 3) subjecting the compound of formula (XII) to oxidation.

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$$R_{6}$$
 R_{7}
 R_{2}
 R_{4}
 R_{1}
 R_{2}
 R_{3}

$$R_8$$
 R_4
 R_8
 R_8
 R_8
 R_8
 R_8

$$R_6$$
 R_7 R_8 R_8 (XII)

wherein,

 R_2 to R_7 have the same meanings as defined in claim 5.

- 8. (As originally filed) The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:
 - 1) reacting compounds of formula (IX) and (XIII) to obtain a compound of formula (XIV); and
 - 2) subjecting the compound of formula (XIV) to cyclization.

wherein,

 R_2 to R_7 have the same meanings as defined in claim 5.

9. (As originally filed) The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:

1) subjecting a compound of formula (XV) to bromination obtain a compound of formula (XVI); and

2) subjecting the compound of formula (XVI) to a carbon-carbon coupling reaction in the presence of a metal catalyst, or to a substitution reaction using a nucleophile.

$$\begin{matrix} R_{0} & \begin{matrix} R_{7} & 0 \\ R_{0} & \begin{matrix} R_{4} \end{matrix} & \begin{matrix} R_{3} \end{matrix} & (XV) \end{matrix}$$

$$R_{8}$$
 R_{4}
 R_{8}
 R_{4}
 R_{8}
 R_{4}
 R_{5}
 R_{4}
 R_{5}

wherein,

R₃ to R₇ have the same meanings as defined in claim 5.

10. (Currently amended) The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:

1) subjecting a compound of formula (XVII) to bromination to obtain a compound of formula (XVIII); and

2) subjecting the compound of formula (XVIII) to a carbon-carbon coupling reaction in the presence of a metal catalyst, or to a substitution reaction using a nucleophile.

wherein,

 R_2 and R_4 to R_7 have the same meanings as defined in claim 5.

11. (As originally filed) The process of claim 5, wherein the indenone compound of formula (II) is prepared by subjecting a compound of formula (XIX) to an acylation reaction, a halogenation reaction followed by a substitution reaction by a nucleophile, or a carbon-carbon coupling reaction in the presence of a metal catalyst.

$$\gamma$$
 -(CH₂)n- R_2

$$R_3$$
(XIX)

wherein,

 R_2 and R_3 have the same meanings as defined in claim 5, Y is hydroxy, thiol, amino C_{1-6} alkyl or halogen, and n is an integer in the range of 0 to 5.

12. (Currently amended) A pharmaceutical composition for modulating the activities of peroxisome proliferator activated receptors (PPARs) comprising a therapeutically effective amount of the compound or salt defined in claim 1 as an active ingredient together with a pharmaceutically acceptable carrier.

13. (Canceled)